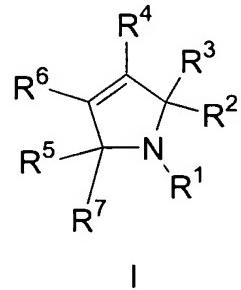


In the claims:

1. (Currently amended) A compound of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof,

wherein:

a is 0 or 1;
b is 0 or 1;
m is 0, 1, or 2;
n is 0;
r is 0 or 1;
s is 0 or 1;

R¹ is selected from:

- 1) (C₁-C₆-alkylene)_n(C=O)C₁-C₁₀ alkyl, and
- 2) (C₁-C₆-alkylene)_n(C=O)aryl,
- 3) (C₁-C₆-alkylene)_n(C=O)C₂-C₁₀ alkenyl,
- 4) (C₁-C₆-alkylene)_n(C=O)C₂-C₁₀ alkynyl,
- 5) (C₁-C₆-alkylene)_n(C=O)C₃-C₈ cycloalkyl,
- 6) (C₁-C₆-alkylene)_n(C=O)heterocyclyl,
- 7) (C₁-C₆-alkylene)_n(C=O)NR^cRC^c,
- 8) (C₁-C₆-alkylene)_nSO₂NR^eRe^e,
- 9) (C₁-C₆-alkylene)_nSO₂C₁-C₁₀ alkyl,
- 10) (C₁-C₆-alkylene)_nSO₂-aryl,
- 11) (C₁-C₆-alkylene)_nSO₂-heterocyclyl,

- 12) $(C_1-C_6\text{-alkylene})_nSO_2-C_3-C_8\text{-cycloalkyl}$,
- 13) $(C_1-C_6\text{-alkylene})_nP(=O)R^dR^{d'}$,
- 14) aryl;
- 15) heterocyclyl;
- 16) $C_1-C_{10}\text{ alkyl}$;
- 17) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_1-C_{10}\text{ alkyl}$,
- 18) $(C_1-C_6\text{-alkylene})_n(C=O)O\text{-aryl}$,
- 19) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_2-C_{10}\text{ alkenyl}$,
- 20) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_2-C_{10}\text{ alkynyl}$,
- 21) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_3-C_8\text{-cycloalkyl}$,
- 22) $(C_1-C_6\text{-alkylene})_n(C=O)O\text{-heterocyclyl}$,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R² and R⁶ are independently selected from:

- 1) aryl,
- 2) $C_1-C_6\text{ aralkyl}$,
- 3) $C_3-C_8\text{-cycloalkyl}$, and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R³ is selected from:

- 1) $C_1-C_{10}\text{ alkyl-O-Rg}$,
- 2) $C_2-C_{10}\text{ alkenyl-O-Rg}$,
- 3) $C_2-C_{10}\text{ alkynyl-O-Rg}$,
- 4) $(C_1-C_6\text{-alkylene})_nC_3-C_8\text{-cycloalkyl-O-Rg}$,
- 5) $C_1-C_{10}\text{ alkyl-(C=O)b-NRfRf'}$,
- 6) $C_2-C_{10}\text{ alkenyl-(C=O)b-NRfRf'}$,
- 7) $C_2-C_{10}\text{ alkynyl-(C=O)b-NRfRf'}$,
- 8) $(C_1-C_6\text{-alkylene})_nC_3-C_8\text{-cycloalkyl-(C=O)b-NRfRf'}$,
- 9) $C_1-C_{10}\text{ alkyl-S(O)m-Rg}$,
- 10) $C_2-C_{10}\text{ alkenyl-S(O)m-Rg}$,

11) ~~C₂-C₁₀ alkynyl S(O)_m R₅~~,

12) ~~(C₁-C₆ alkylene)_n C₃-C₈ cycloalkyl S(O)_m R₅~~,

said alkyl, alkenyl, alkynyl and cycloalkyl are is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ is selected from:

1) H, and

2) C₁-C₁₀ alkyl,

3) aryl,

4) C₂-C₁₀ alkenyl,

5) C₂-C₁₀ alkynyl,

6) C₁-C₆ perfluoroalkyl,

7) C₁-C₆ aralkyl,

8) C₃-C₈ cycloalkyl, and

9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R⁵ and R⁷ are independently selected from:

1) H, and

2) C₁-C₁₀ alkyl,

3) aryl,

4) C₂-C₁₀ alkenyl,

5) C₂-C₁₀ alkynyl,

6) C₁-C₆ perfluoroalkyl,

7) C₁-C₆ aralkyl,

8) C₃-C₈ cycloalkyl, and

9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰; or

R⁵ and R⁷ are combined to form an oxo or a sulfoxide;

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,
- 10) (C=O)_rO_s(C₀-C₆)alkylene-aryl,
- 11) (C=O)_rO_s(C₀-C₆)alkylene-heterocyclyl,
- 12) (C=O)_rO_s(C₀-C₆)alkylene-N(R^b)₂,

- 13) C(O)R^a,
- 14) (C₀-C₆)alkylene-CO₂R^a,
- 15) C(O)H,
- 16) (C₀-C₆)alkylene-CO₂H, and
- 17) C(O)N(R^b)₂,
- 18) S(O)_mR^a, and
- 19) S(O)₂N(R^b)₂;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R^b, OH, (C₁-C₆)alkoxy, halogen, CO₂H, CN, O(C=O)C₁-C₆ alkyl, oxo, NO₂ and N(R^b)₂;

R¹² and R¹³ are independently selected from:

- 1) H,
- 2) (C=O)O_bC₁-C₁₀ alkyl,
- 3) (C=O)O_bC₃-C₈ cycloalkyl,
- 4) (C=O)O_baryl,
- 5) (C=O)O_bheterocyclyl,
- 6) C₁-C₁₀ alkyl,
- 7) aryl,
- 8) C₂-C₁₀ alkenyl,
- 9) C₂-C₁₀ alkynyl,
- 10) heterocyclyl,
- 11) C₃-C₈ cycloalkyl,
- 12) SO₂R^a, and
- 13) (C=O)NR^b₂,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R¹¹, or

R¹² and R¹³ can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fRF' or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^{c'} are independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or

R^c and R^{c'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^{d'} are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^{d'} can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹; and

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

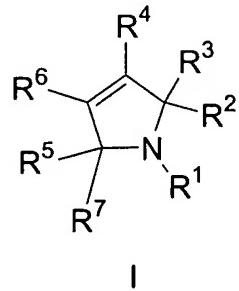
R^f and R^{f'} are independently selected from: H, (C₁-C₆)alkyl, aryl, NH₂, OH, OR^a, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fRF', S(O)₂R^a and -(C₁-C₆)alkyl-N(R^b)₂, wherein the alkyl is optionally substituted with one, two or three substituents selected from R¹¹; or

R^f and R^{f'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R_g is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

2. (Currently amended) The compound according to Claim 1 of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof,

wherein:

- a is 0 or 1;
b is 0 or 1;
m is 0, 1, or 2;
n is 0 or 1;
r is 0 or 1;
s is 0 or 1;

R¹ is selected from:

- 1) (C₁-C₆-alkylene)_n(C=O)C₁-C₁₀ alkyl,
- 2) (C₁-C₆-alkylene)_n(C=O)aryl,
- 3) (C₁-C₆-alkylene)_n(C=O)C₂-C₁₀ alkenyl,
- 4) (C₁-C₆-alkylene)_n(C=O)C₂-C₁₀ alkynyl,
- 5) (C₁-C₆-alkylene)_n(C=O)C₃-C₈ cycloalkyl,
- 6) (C₁-C₆-alkylene)_n(C=O)heterocyclyl,
- 2) 7) (C₁-C₆-alkylene)_n(C=O)NR^cRC',

- 7) ~~(C₁-C₆-alkylene)_nSO₂NR_eR_{e'},~~
- 8) ~~(C₁-C₆-alkylene)_nSO₂C₁-C₁₀-alkyl,~~
- 9) ~~(C₁-C₆-alkylene)_nSO₂-aryl,~~
- 10) ~~(C₁-C₆-alkylene)_nSO₂-heterocyclyl,~~
- 11) ~~(C₁-C₆-alkylene)_nSO₂-C₃-C₈-cycloalkyl,~~
- 12) ~~(C₁-C₆-alkylene)_nP(=O)R_dR_{d'},~~
- 13) ~~aryl;~~
- 14) ~~heterocyclyl;~~
- 15) ~~C₁-C₁₀-alkyl;~~
- 16) ~~(C₁-C₆-alkylene)_n(C=O)O-C₁-C₁₀-alkyl,~~
- 17) ~~(C₁-C₆-alkylene)_n(C=O)O-aryl,~~
- 18) ~~(C₁-C₆-alkylene)_n(C=O)O-C₂-C₁₀-alkenyl,~~
- 19) ~~(C₁-C₆-alkylene)_n(C=O)O-C₂-C₁₀-alkynyl,~~
- 20) ~~(C₁-C₆-alkylene)_n(C=O)O-C₃-C₈-cycloalkyl,~~
- 21) ~~(C₁-C₆-alkylene)_n(C=O)O-heterocyclyl,~~

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R² and R⁶ are independently selected from:

- 1) aryl,
- 2) C₁-C₆-aralkyl,
- 3) C₃-C₈-cycloalkyl, and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R³ is selected from:

- 1) C₁-C₁₀ alkyl-O-Rg,
- 2) C₂-C₁₀ alkenyl-O-Rg,
- 3) C₂-C₁₀ alkynyl-O-Rg,
- 4) (C₁-C₆-alkylene)_nC₃-C₈-cycloalkyl-O-Rg,
- 5) C₁-C₁₀ alkyl-(C=O)_b-NR^fR^{f'},
- 6) C₂-C₁₀ alkenyl-(C=O)_bNR^fR^{f'},
- 7) C₂-C₁₀ alkynyl-(C=O)_bNR^fR^{f'},

8) ~~(C₁-C₆ alkylene)_nC₃-C₈ cycloalkyl (C=O)_bNR^fR^{f'},~~

9) ~~C₁-C₁₀ alkyl S(O)_m-R₈,~~

10) ~~C₂-C₁₀ alkenyl S(O)_m-R₈,~~

11) ~~C₂-C₁₀ alkynyl S(O)_m-R₈,~~

12) ~~(C₁-C₆ alkylene)_nC₃-C₈ cycloalkyl S(O)_m-R₈,~~

said alkyl, alkenyl, alkynyl and cycloalkyl are is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ is selected from:

1) H,

2) C₁-C₁₀ alkyl,

3) aryl,

4) C₂-C₁₀ alkenyl,

5) C₂-C₁₀ alkynyl,

6) C₁-C₆ perfluoroalkyl,

7) C₄-C₆ aralkyl,

8) C₃-C₈ cycloalkyl, and

9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R⁵ and R⁷ are independently selected from:

1) H,

2) C₁-C₁₀ alkyl,

3) aryl,

4) C₂-C₁₀ alkenyl,

5) C₂-C₁₀ alkynyl,

6) C₁-C₆ perfluoroalkyl,

7) C₄-C₆ aralkyl,

8) C₃-C₈ cycloalkyl, and

9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰; or

~~R⁵ and R⁷ are combined to form an exo or a sulfoxo;~~

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,

- 10) $(C=O)_r Os(C_0-C_6)$ alkylene-aryl,
- 11) $(C=O)_r Os(C_0-C_6)$ alkylene-heterocyclyl,
- 12) $(C=O)_r Os(C_0-C_6)$ alkylene-N(R^b)₂,
- 13) C(O)R^a,
- 14) (C_0-C_6) alkylene-CO₂R^a,
- 15) C(O)H,
- 16) (C_0-C_6) alkylene-CO₂H, and
- 17) C(O)N(R^b)₂,
- 18) S(O)_mR^a, and
- 19) S(O)₂N(R^b)₂;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R^b, OH, (C₁-C₆)alkoxy, halogen, CO₂H, CN, O(C=O)C₁-C₆ alkyl, oxo, and N(R^b)₂;

R¹² and R¹³ are independently selected from:

- 1) H,
- 2) (C=O)O_bC₁-C₁₀ alkyl,
- 3) (C=O)O_bC₃-C₈ cycloalkyl,
- 4) (C=O)O_baryl,
- 5) (C=O)O_bheterocyclyl,
- 6) C₁-C₁₀ alkyl,
- 7) aryl,
- 8) C₂-C₁₀ alkenyl,
- 9) C₂-C₁₀ alkynyl,
- 10) heterocyclyl,
- 11) C₃-C₈ cycloalkyl,
- 12) SO₂R^a, and
- 13) (C=O)NR^b₂,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R¹¹, or

R¹² and R¹³ can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in

addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^f' or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^{c'} are independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or R^c and R^{c'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^{d'} are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^{d'} can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from N^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹; and

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

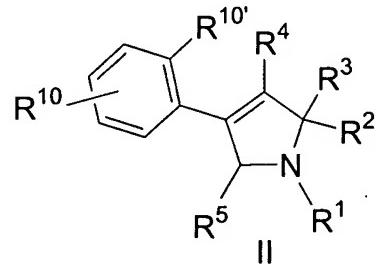
R^f and R^{f'} are independently selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂, or

R^f and R^{f'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R_g is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

3. (Currently amended) The compound according to Claim 2 of Formula II:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;

b is 0 or 1;

m is 0, 1, or 2;

r is 0 or 1;

s is 0 or 1;

R¹ is selected from:

- 1) (C=O)C₁-C₁₀ alkyl, and
- 2) (C=O)aryl,
- 3) (C=O)C₂-C₁₀ alkenyl,
- 4) (C=O)C₂-C₁₀ alkynyl,
- 5) (C=O)C₃-C₈ cycloalkyl,
- 6) (C=O)heterocyclyl,
- 7) (C=O)NR^cRC^c,
- 8) SO₂NR^eRE^e,
- 9) SO₂C₁-C₁₀ alkyl,
- 10) SO₂ aryl,
- 11) SO₂ heterocyclyl,

12) $\text{SO}_2\text{C}_3\text{-}\text{C}_8\text{-cycloalkyl}$, and

13) $\text{P}(\text{-O})\text{R}^d\text{R}^{d'}$,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R² is selected from:

1) aryl,

2) $\text{C}_1\text{-}\text{C}_6$ aralkyl,

3) $\text{C}_3\text{-}\text{C}_8$ cycloalkyl, and

4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R³ is selected from:

1) C₁-C₁₀ alkyl-O-Rg,

2) $\text{C}_3\text{-}\text{C}_8$ cycloalkyl-O-Rg,

2 3) C₁-C₁₀ alkyl- NR^fR^{f'},

4) $\text{C}_3\text{-}\text{C}_8$ cycloalkyl- NR^fR^{f'},

said alkyl and cycloalkyl are is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ and R⁵ are independently selected from:

1) H, and

2) C₁-C₁₀ alkyl,

3) aryl,

4) C₂-C₁₀ alkenyl,

5) C₂-C₁₀ alkynyl,

6) C₁-C₆ perfluoroalkyl,

7) C₁-C₆ aralkyl,

8) $\text{C}_3\text{-}\text{C}_8$ cycloalkyl, and

9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R^{10'} is halogen;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,
- 10) (C=O)_rO_s(C₀-C₆)alkylene-aryl,

- 11) $(C=O)_r Os(C_0-C_6)$ alkylene-heterocyclyl,
- 12) $(C=O)_r Os(C_0-C_6)$ alkylene-N(R^b)₂,
- 13) C(O)R^a,
- 14) (C_0-C_6) alkylene-CO₂R^a,
- 15) C(O)H,
- 16) (C_0-C_6) alkylene-CO₂H, and
- 17) C(O)N(R^b)₂,
- 18) S(O)_mR^a, and
- 19) S(O)₂N(R^b)₂;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R^b, OH, (C₁-C₆)alkoxy, halogen, CO₂H, CN, O(C=O)C₁-C₆ alkyl, oxo, and N(R^b)₂;

R¹² and R¹³ are independently selected from:

- 1) H,
- 2) (C=O)O_bC₁-C₁₀ alkyl,
- 3) (C=O)O_bC₃-C₈ cycloalkyl,
- 4) (C=O)Obaryl,
- 5) (C=O)Obheterocyclyl,
- 6) C₁-C₁₀ alkyl,
- 7) aryl,
- 8) C₂-C₁₀ alkenyl,
- 9) C₂-C₁₀ alkynyl,
- 10) heterocyclyl,
- 11) C₃-C₈ cycloalkyl,
- 12) SO₂R^a, and
- 13) (C=O)NR^b₂,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R¹¹, or

R¹² and R¹³ can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^f' or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^c' are independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or

R^c and R^c' can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^d' are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^d' can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^f and R^f' are independently selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂, or

R^f and R^f' can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

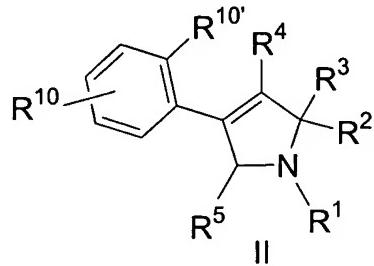
R_g is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

4. (Cancelled)

5. (Canceled)

6 (Original) The compound according to Claim 3 of the Formula II, or a pharmaceutically acceptable salt or stereoisomer thereof, wherein R² is phenyl, optionally substituted with one or two substituents selected from R¹⁰.

7. (Currently amended) The compound according to Claim 1 of the formula II:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;

b is 0 or 1;

m is 0, 1, or 2;

r is 0 or 1;

s is 0 or 1;

R¹ is selected from:

- 1) (C=O)C₁-C₁₀ alkyl, and
- 2) (C=O)aryl,

- 3) $(C=O)C_2-C_{10}$ alkenyl,
- 4) $(C=O)C_2-C_{10}$ alkynyl,
- 5) $(C=O)C_3-C_8$ cycloalkyl,
- 6) $(C=O)$ heterocyclyl,
- 2) 7) $(C=O)NR^cR^{c'}$,
- 7) $SO_2NR^eR^{e'}$,
- 8) $SO_2C_1-C_{10}$ alkyl,
- 9) SO_2 aryl,
- 10) SO_2 heterocyclyl,
- 11) $SO_2C_3-C_8$ cycloalkyl, and
- 12) $P(=O)R^dR^{d'}$,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R² is phenyl, optionally substituted with one or more substituents selected from R¹⁰;

R³ is selected from:

- 1) C₁-C₁₀ alkyl-O-R_g,
- 2) C₁-C₁₀ alkyl- NR^fR^{f'},

said alkyl is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ and R⁵ are independently selected from:

- 1) H, and
- 2) C₁-C₁₀ alkyl,
- 3) aryl,
- 4) C₂-C₁₀ alkenyl,
- 5) C₂-C₁₀ alkynyl,
- 6) C₁-C₆ perfluoroalkyl,
- 7) C₁-C₆ aralkyl,
- 8) C₃-C₈ cycloalkyl, and
- 9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R^{10'} is halogen;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,
- 10) (C=O)_rO_s(C₀-C₆)alkylene-aryl,

- 11) $(C=O)_r O_s (C_0-C_6)$ alkylene-heterocyclyl,
- 12) $(C=O)_r O_s (C_0-C_6)$ alkylene-N(R^b)₂,
- 13) C(O)R^a,
- 14) (C_0-C_6) alkylene-CO₂R^a,
- 15) C(O)H,
- 16) (C_0-C_6) alkylene-CO₂H, and
- 17) C(O)N(R^b)₂,
- 18) S(O)_mR^a, and
- 19) S(O)₂N(R^b)₂;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R^b, OH, (C₁-C₆)alkoxy, halogen, CO₂H, CN, O(C=O)C₁-C₆ alkyl, oxo, and N(R^b)₂;

R¹² and R¹³ are independently selected from:

- 1) H,
- 2) (C=O)O_bC₁-C₁₀ alkyl,
- 3) (C=O)O_bC₃-C₈ cycloalkyl,
- 4) (C=O)O_baryl,
- 5) (C=O)O_bheterocyclyl,
- 6) C₁-C₁₀ alkyl,
- 7) aryl,
- 8) C₂-C₁₀ alkenyl,
- 9) C₂-C₁₀ alkynyl,
- 10) heterocyclyl,
- 11) C₃-C₈ cycloalkyl,
- 12) SO₂R^a, and
- 13) (C=O)NR^b₂,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R¹¹, or

R¹² and R¹³ can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^f' or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^c' are independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or

R^c and R^c' can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^d' are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^d' can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^f and R^f' are independently selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂, or

R^f and R^f' can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in

addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R_g is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

8. (Amended) A compound selected from:

4-(2,5-Difluorophenyl)-2-(hydroxymethyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-(hydroxymethyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-(methoxymethyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-[(2-hydroxyethoxy)methyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[(2-Aminoethoxy)methyl]-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-({[2-(dimethylamino)ethyl]amino}methyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

3-{4-(2,5-Difluorophenyl)-1-[(dimethylamino)carbonyl]-2-phenyl-2,5-dihydro-1H-pyrrol-2-yl}prop-2-en-1-aminium amine;

2-(3-Hydroxypropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-Aminopropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-[3-(dimethylamino)propyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(1-hydroxyethyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

[4-(2,5-difluorophenyl)-2-phenyl-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1*H*-pyrrol-2-yl]methanol;

2-({[tert-butyl(dimethyl)silyl]oxy}methyl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-*N*-[1-(*N,N*-dimethylglycyl)piperidin-4-yl]-2-(hydroxymethyl)-*N*-methyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(hydroxymethyl)-*N*-methyl-*N*-[1-(morpholin-4-ylacetyl)piperidin-4-yl]-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(hydroxymethyl)-*N*-methyl-2-phenyl-*N*-piperidin-3-yl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

N-[1-(2,2-difluoroethyl)piperidin-4-yl]-4-(2,5-difluorophenyl)-2-(hydroxymethyl)-*N*-methyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-*N*-[1-(2-hydroxyethyl)piperidin-4-yl]-2-(hydroxymethyl)-*N*-methyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

(2*S*)-4-(2,5-difluorophenyl)-*N*-[1-(2-fluoroethyl)piperidin-4-yl]-2-(hydroxymethyl)-*N*-methyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(hydroxymethyl)-*N*-methyl-*N*-{1-[(methylsulfonyl)methyl]piperidin-4-yl}-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-*N*-{1-[2-fluoro-1-(fluoromethyl)ethyl]piperidin-4-yl}-2-(hydroxymethyl)-*N*-methyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

(2*S*)-*N*-(1-cyclopropylpiperidin-4-yl)-4-(2,5-difluorophenyl)-2-(hydroxymethyl)-*N*-methyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

benzyl {4-[{[4-(2,5-difluorophenyl)-2-(hydroxymethyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}(methyl)amino]piperidin-1-yl}acetate;

{4-[{[4-(2,5-difluorophenyl)-2-(hydroxymethyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}(methyl)amino]piperidin-1-yl}acetic acid;

methyl {4-[{[4-(2,5-difluorophenyl)-2-(hydroxymethyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}(methyl)amino]piperidin-1-yl}acetate;

4-(2,5-difluorophenyl)-2-(methoxymethyl)-*N*-methyl-*N*-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxypropyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-{3-[(2,2-difluoroethyl)amino]propyl}-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-{3-[(2,2-difluoroethyl)(methyl)amino]propyl}-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-(3-aminopropyl)-4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-[3-(acetylamino)propyl]-4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-{3-[(methylsulfonyl)amino]propyl}-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

methyl 3-{4-(5-chloro-2-fluorophenyl)-1-[(dimethylamino)carbonyl]-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl}propylcarbamate;

2-{3-[(aminocarbonyl)amino]propyl}-4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

3-{4-(2,5-difluorophenyl)-1-[(dimethylamino)carbonyl]-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl}propanoic acid;

2-(3-anilino-3-oxopropyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydrazino-3-oxopropyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-[3-(hydroxyamino)-3-oxopropyl]-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-(2,2-difluoro-3-hydroxypropyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-(3-amino-2,2-difluoropropyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-[3-(dimethylamino)propyl]-*N*-methyl-2-phenyl-*N*-tetrahydro-2*H*-pyran-4-yl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

1-{4-(2,5-difluorophenyl)-2-[3-(dimethylamino)propyl]-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl}-2-methyl-1-oxopropan-2-ol;

3-[(2*S*)-1-[(2*S*)-2-amino-2-cyclopropylethanoyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl]-*N,N*-dimethylpropan-1-amine; and
(2*S*)-2-(3-amino-4,4-difluorobutyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

or a pharmaceutically acceptable salt or stereoisomer thereof.

9. (Original) The compound according to Claim 8 which is selected from:

4-(2,5-Difluorophenyl)-2-(hydroxymethyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[(2-Aminoethoxy)methyl]-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-Aminopropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-[3-(dimethylamino)propyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

or a pharmaceutically acceptable salt or stereoisomer thereof.

10. (Canceled)

11. (Original) The compound according to Claim 1 selected from:

2-[(2-Aminoethoxy)methyl]-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide hydrochloride salt;

4-(2,5-Difluorophenyl)-2-({[2-(dimethylamino)ethyl]amino}methyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide bis TFA salt;

2-(3-Aminopropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide hydrochloride salt;

4-(2,5-Difluorophenyl)-2-[3-(dimethylamino)propyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide TFA salt;

4-(2,5-difluorophenyl)-N-[1-(glycyl)piperidin-4-yl]-2-(hydroxymethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide TFA salt;

3-[(2S)-1-[(2S)-2-amino-2-cyclopropylethanoyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl]-*N,N*-dimethylpropan-1-amine bis-TFA salt;

3-[(2R)-1-[(2S)-2-amino-2-cyclopropylethanoyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl]-*N,N*-dimethylpropan-1-amine bis-TFA salt;

4-(2,5-difluorophenyl)-2-[3-(dimethylamino)propyl]-*N*-methyl-*N*-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide bis-TFA salt;

4-(2,5-difluorophenyl)-2-[3-(ethylamino)propyl]-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide TFA salt;

4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2-{3-[(pyridin-4-ylmethyl)amino]propyl}-2,5-dihydro-1*H*-pyrrole-1-carboxamide bis-TFA salt; and

4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-(3-{{[(4-methyl-1*H*-imidazol-2-yl)methyl]amino}propyl}-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide bis-TFA salt.

12. (Original) A pharmaceutical composition that is comprised of a compound in accordance with Claim 1 and a pharmaceutically acceptable carrier.

13. (Cancelled)

14. (Cancelled)

15. (Cancelled)

16. (Canceled)

17. (Canceled)

18. (Canceled)

19. (Canceled)
20. (Canceled)
21. (Canceled)
22. (Canceled)
23. (Canceled)
24. (Canceled)
25. (Cancelled)
26. (Cancelled)
27. (Canceled)
28. (Canceled)
29. (Canceled)
30. (Canceled)
31. (Canceled)
32. (Canceled)
33. (Canceled)
34. (Canceled)
35. (Canceled)

36. (Canceled)

37. (Canceled)

38. (Cancelled)